

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101220\
 Data File : BM027652.D
 Acq On : 12 Oct 2020 10:11
 Operator : JU/CG
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4008

Quant Time: Oct 12 10:51:31 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM100820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 12 10:50:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	605	0.40	ng/ul	0.00
2) Naphthalene-d8	10.57	136	1769	0.40	ng/ul	0.00
7) Acenaphthene-d10	14.41	164	1237	0.40	ng/ul	0.00
11) Phenanthrene-d10	17.16	188	3210	0.40	ng/ul	0.00
17) Chrysene-d12	21.33	240	3381	0.40	ng/ul	0.00
21) Perylene-d12	23.59	264	3529	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.16	152	1345	0.41	ng/ul	0.00
15) Fluoranthene-d10	19.18	212	4241	0.43	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.62	128	2052	0.407	ng/ul	97
5) 2-Methylnaphthalene	12.23	142	1519	0.416	ng/ul	100
6) 1-Methylnaphthalene	12.45	142	1726	0.453	ng/ul	100
8) Acenaphthylene	14.13	152	2213	0.412	ng/ul	97
9) Acenaphthene	14.48	153	1763	0.415	ng/ul	96
10) Fluorene	15.46	166	2279	0.421	ng/ul	98
12) Pentachlorophenol	16.80	266	549	0.421	ng/ul	95
13) Phenanthrene	17.20	178	3995	0.411	ng/ul	97
14) Anthracene	17.29	178	3764	0.414	ng/ul	98
16) Fluoranthene	19.21	202	5352	0.436	ng/ul	97
18) Pyrene	19.57	202	5427	0.463	ng/ul#	97
19) Benzo(a)anthracene	21.32	228	5364	0.410	ng/ul	98
20) Chrysene	21.37	228	5698	0.410	ng/ul	99
22) Benzo(b)fluoranthene	22.91	252	6044	0.403	ng/ul	97
23) Benzo(k)fluoranthene	22.95	252	6061	0.404	ng/ul	97
24) Benzo(a)pyrene	23.49	252	5396	0.411	ng/ul#	95
25) Indeno(1,2,3-cd)pyrene	25.85	276	7046	0.399	ng/ul#	100
26) Dibenzo(a,h)anthracene	25.87	278	5879	0.403	ng/ul	98
27) Benzo(a,h,i)perylene	26.55	276	6045	0.391	ng/ul	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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